

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041817\
 Data File : BG026653.D
 Acq On : 18 Apr 2017 17:43
 Operator : SJ/MA
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG041817

Quant Time: Apr 18 18:56:13 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 18 18:04:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	162594	20.00	ng	-0.02
21) Naphthalene-d8	10.83	136	698922	20.00	ng	-0.02
38) Acenaphthene-d10	14.64	164	432424	20.00	ng	-0.02
63) Phenanthrene-d10	17.37	188	1092441	20.00	ng	-0.02
75) Chrysene-d12	21.61	240	1227431	20.00	ng	-0.02
86) Perylene-d12	24.78	264	1232870	20.00	ng	-0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	729103	81.52	ng	-0.02
7) Phenol-d6	7.19	99	976267	83.30	ng	-0.02
23) Nitrobenzene-d5	9.19	82	825817	82.45	ng	-0.02
41) 2,4,6-Tribromophenol	16.12	330	497717	82.60	ng	-0.02
44) 2-Fluorobiphenyl	13.26	172	2297821	79.14	ng	-0.02
78) Terphenyl-d14	19.97	244	3311087	79.64	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	3.48	88	188145	40.23	ng	# 71
3) Pyridine	3.89	79	437772	41.81	ng	# 62
4) n-Nitrosodimethylamine	3.80	42	123634	40.59	ng	# 1
6) Aniline	7.35	93	629813	43.03	ng	# 87
8) 2-Chlorophenol	7.59	128	441604	41.38	ng	# 79
9) Benzaldehyde	7.16	77	327253	41.22	ng	# 65
10) Phenol	7.22	94	530564	41.61	ng	# 69
11) bis(2-Chloroethyl)ether	7.44	93	412237	40.61	ng	# 80
12) 1,3-Dichlorobenzene	7.92	146	518281	41.21	ng	# 84
13) 1,4-Dichlorobenzene	8.06	146	519612m	40.67	ng	
14) 1,2-Dichlorobenzene	8.38	146	499204	41.35	ng	# 90
15) Benzyl Alcohol	8.26	79	308998	42.16	ng	# 75
16) 2,2'-oxybis(1-Chloropropan	8.56	45	389224	39.63	ng	79
17) 2-Methylphenol	8.46	107	374537	41.92	ng	# 80
18) Hexachloroethane	9.11	117	164861	40.89	ng	94
19) n-Nitroso-di-n-propylamine	8.83	70	298870	41.14	ng	# 70
20) 3+4-Methylphenols	8.79	107	520423	42.39	ng	87
22) Acetophenone	8.85	105	645443	40.41	ng	# 73
24) Nitrobenzene	9.23	77	434194	40.94	ng	# 71
25) Isophorone	9.75	82	847245	41.26	ng	# 86
26) 2-Nitrophenol	9.94	139	268407	41.77	ng	# 62
27) 2,4-Dimethylphenol	10.00	122	429618	41.42	ng	86
28) bis(2-Chloroethoxy)methane	10.23	93	555681	41.40	ng	# 97
29) 2,4-Dichlorophenol	10.48	162	461703	42.56	ng	93
30) 1,2,4-Trichlorobenzene	10.69	180	489748	40.50	ng	99
31) Naphthalene	10.88	128	1403962	40.93	ng	99
32) Benzoic acid	10.13	122	289463	41.33	ng	# 70
33) 4-Chloroaniline	10.99	127	612111	41.60	ng	# 79
34) Hexachlorobutadiene	11.17	225	284399	41.13	ng	99
35) Caprolactam	11.75	113	178890	42.47	ng	# 47
36) 4-Chloro-3-methylphenol	12.10	107	457878	41.91	ng	# 70
37) 2-Methylnaphthalene	12.47	142	1091308	41.34	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	536057	40.25	ng	# 97
40) Hexachlorocyclopentadiene	12.82	237	274203	40.74	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.08	196	384803	40.57	nq	95
43) 2,4,5-Trichlorophenol	13.15	196	404488	40.33	nq	# 91
45) 1,1'-Biphenyl	13.48	154	1400281	39.85	nq	99
46) 2-Chloronaphthalene	13.52	162	1065933	40.03	nq	97
47) 2-Nitroaniline	13.72	65	282325	41.51	nq	# 49
48) Acenaphthylene	14.36	152	1719947	40.17	nq	99
49) Dimethylphthalate	14.09	163	1394529	40.74	nq	98
50) 2,6-Dinitrotoluene	14.20	165	339937	42.47	nq	# 57
51) Acenaphthene	14.70	154	1096792	40.24	nq	97
52) 3-Nitroaniline	14.54	138	364075	41.70	nq	# 64
53) 2,4-Dinitrophenol	14.75	184	164085	39.49	nq	# 77
54) Dibenzofuran	15.03	168	1722771	40.56	nq	# 89
55) 4-Nitrophenol	14.85	139	314355	42.65	nq	# 38
56) 2,4-Dinitrotoluene	14.99	165	490077	43.11	nq	# 68
57) Fluorene	15.68	166	1420306	40.61	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.26	232	407399	39.94	nq	# 95
59) Diethylphthalate	15.44	149	1413143	40.61	nq	99
60) 4-Chlorophenyl-phenylether	15.67	204	717237	41.09	nq	# 87
61) 4-Nitroaniline	15.70	138	398230	42.77	nq	# 49
62) Azobenzene	15.96	77	1097011	40.85	nq	79
64) 4,6-Dinitro-2-methylphenol	15.76	198	313120	42.44	nq	86
65) n-Nitrosodiphenylamine	15.88	169	1275794	40.48	nq	99
66) 4-Bromophenyl-phenylether	16.56	248	477726	40.02	nq	94
67) Hexachlorobenzene	16.68	284	528097	40.24	nq	# 88
68) Atrazine	16.82	200	504978	40.98	nq	98
69) Pentachlorophenol	17.03	266	329719	41.59	nq	97
70) Phenanthrene	17.41	178	2270294	40.36	nq	98
71) Anthracene	17.50	178	2349872	40.43	nq	97
72) Carbazole	17.77	167	2248717	40.40	nq	97
73) Di-n-butylphthalate	18.32	149	2505272	40.14	nq	# 95
74) Fluoranthene	19.42	202	2690283	41.25	nq	97
76) Benzidine	19.59	184	1419461	43.07	nq	98
77) Pyrene	19.77	202	2724462	41.23	nq	97
79) Butylbenzylphthalate	20.65	149	1167734	41.07	nq	# 81
80) Benzo(a)anthracene	21.60	228	2660170	40.13	nq	99
81) 3,3'-Dichlorobenzidine	21.51	252	1110830	40.40	nq	100
82) Chrysene	21.66	228	2459328	40.32	nq	98
83) Bis(2-ethylhexyl)phthalate	21.49	149	1658955	40.49	nq	96
84) Di-n-octyl phthalate	22.68	149	2816949	41.17	nq	# 81
85) Indeno(1,2,3-cd)pyrene	28.40	276	3429072	41.00	nq	# 88
87) Benzo(b)fluoranthene	23.78	252	2888395	40.92	nq	# 96
88) Benzo(k)fluoranthene	23.84	252	2733155	40.17	nq	# 95
89) Benzo(a)pyrene	24.63	252	2767464	40.61	nq	# 96
90) Dibenzo(a,h)anthracene	28.45	278	2891213	41.30	nq	# 93
91) Benzo(g,h,i)perylene	29.52	276	2850590	40.72	nq	# 89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

